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AND CURRENT DISCOVER FILE IS DATED 19 SEPTEMBER 2007.

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NEWS LOGIN Welcome Banner and News Items
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STRUCTURE FILE UPDATES: 27 JAN 2008 HIGHEST RN 1000849-38-6
DICTIONARY FILE UPDATES: 27 JAN 2008 HIGHEST RN 1000849-38-6

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TSCA INFORMATION NOW CURRENT THROUGH June 29, 2007

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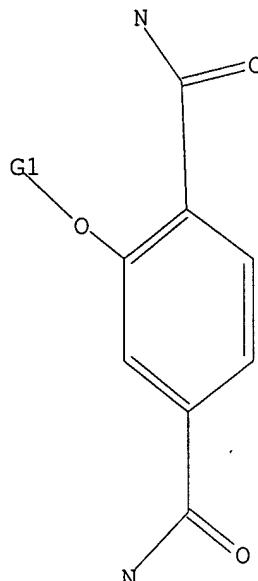
REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stnqgen/stndoc/properties.html>

=> Uploading C:\Documents and Settings\jcho2\My Documents\10588478-a.str

L1 STRUCTURE UPLOADED

=> d 11
L1 HAS NO ANSWERS
L1 STR



G1 Ak,Cy

Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SEARCH INITIATED 17:26:07 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 173 TO ITERATE

100.0% PROCESSED 173 ITERATIONS 22 ANSWERS
SEARCH TIME: 00.00.01 .

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
                        BATCH **COMPLETE**
PROJECTED ITERATIONS: 2671 TO 4249
PROJECTED ANSWERS: 159 TO 721
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12 22 SEA SSS SAM 1.1

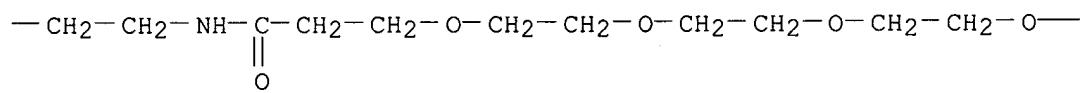
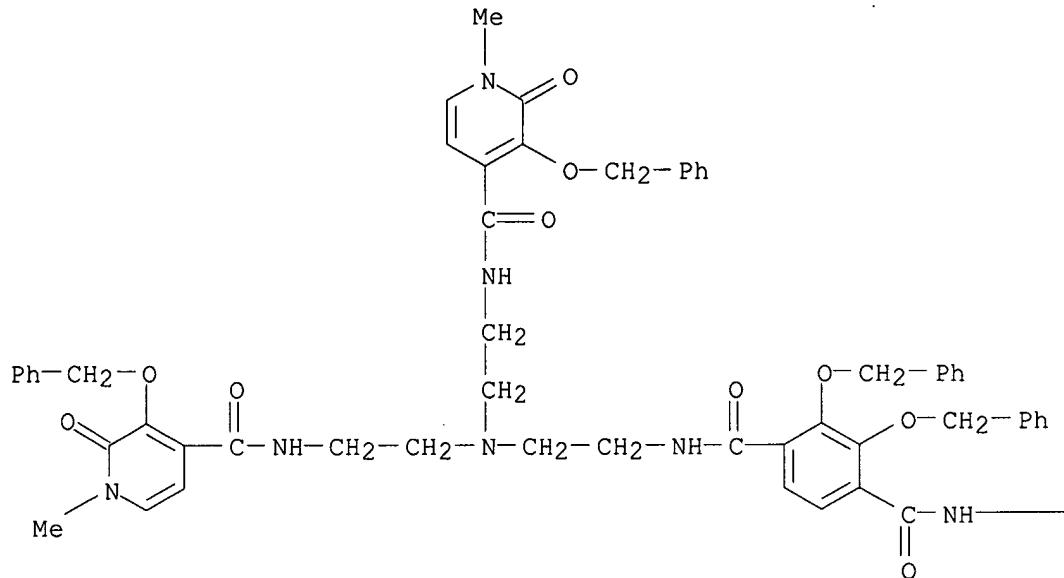
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FULL SCREEN SEARCH COMPLETED - 3389 TO ITERATE
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100.0% PROCESSED 3389 ITERATIONS 564 ANSWERS
SEARCH TIME: 00.00.01

L3 564 SEA SSS FUL L1

=> d scan

L3 564 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
 IN 1,4-Benzenedicarboxamide, N-[2-[bis[2-[[[1,2-dihydro-1-methyl-2-oxo-3-
 (phenylmethoxy)-4-pyridinyl]carbonyl]amino]ethyl]amino]ethyl]-N'-(4-oxo-
 7,10,13,16,19,22,25,28,31,34,37,40-dodecaoxa-3-azahentetracont-1-yl)-2,3-
 bis(phenylmethoxy)- (9CI)
 MF C84 H112 N8 O23



PAGE 1-C

— CH₂— CH₂— O— CH₂— CH₂— O— CH₂— CH₂— O— CH₂— CH₂— O— CH₂— CH₂— O—

PAGE 1-D

— CH₂— CH₂— O— CH₂— CH₂— O— CH₂— CH₂— OMe

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

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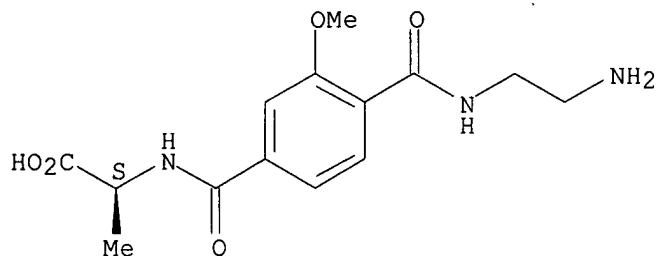
L4 STRUCTURE UPLOADED

=> d 12

L2 ANSWER 1 OF 22 REGISTRY COPYRIGHT 2008 ACS on STN
RN 919772-33-1 REGISTRY
ED Entered STN: 07 Feb 2007
CN L-Alanine, N-[4-[(2-aminoethyl)amino]carbonyl]-3-methoxybenzoyl]-(CA

INDEX NAME)
FS STEREOSEARCH
MF C14 H19 N3 O5
SR CA
LC STN Files: CA, CAPLUS, USPATFULL

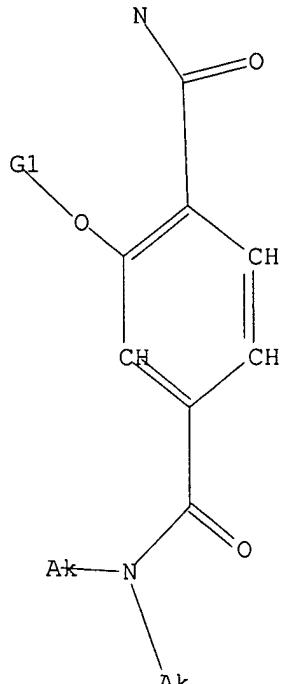
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

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L4 HAS NO ANSWERS
L4 STR



G1 Ak,Cy

Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SEARCH INITIATED 17:31:42 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 173 TO ITERATE

100.0% PROCESSED 173 ITERATIONS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 2671 TO 4249
PROJECTED ANSWERS: 0 TO 0

L5 0 SEA SSS SAM L4

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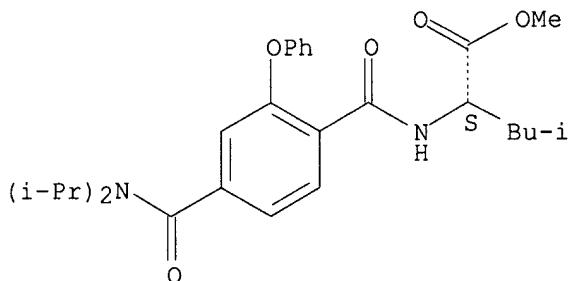
100.0% PROCESSED 3389 ITERATIONS 27 ANSWERS
SEARCH TIME: 00.00.01

L6 27 SEA SSS FUL L4

=> d scan

L6 27 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-phenoxybenzoyl]-,
methyl ester
MF C27 H36 N2 O5

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> file caplus
COST IN U.S. DOLLARS SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST 362.86 363.07

FILE 'CAPLUS' ENTERED AT 17:32:05 ON 28 JAN 2008
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FILE COVERS 1907 - 28 Jan 2008 VOL 148 ISS 5
FILE LAST UPDATED: 27 Jan 2008 (20080127/ED)

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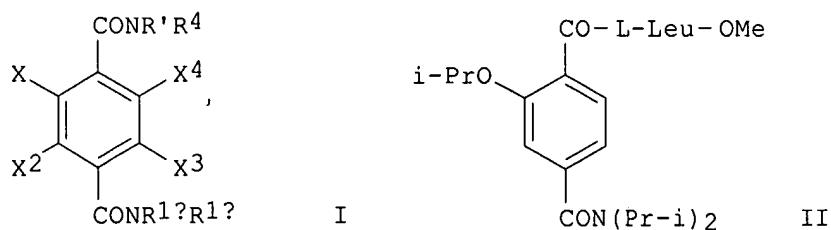
<http://www.cas.org/infopolicy.html>

=> s 16
L7 4 L6

=> d 17 1-4 bib abs hitstr

L7 ANSWER 1 OF 4 CAPLUS COPYRIGHT 2008 ACS on STN
AN 2005:961964 CAPLUS
DN 143:248664
TI Preparation of terephthalamide peptidomimetic compounds for therapeutic use
IN Hamilton, Andrew D.; Yin, Hang
PA Yale University, USA
SO PCT Int. Appl., 48 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1
PATENT NO. KIND DATE APPLICATION NO. DATE

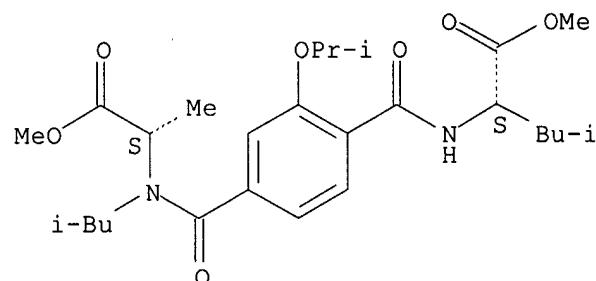
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WO 2005079541 A3 20051103
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CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD,
GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,
LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI,
NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY,
TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW
RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM,
AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK,
EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT,
RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML,
MR, NE, SN, TD, TG
AU 2005215051 A1 20050901 AU 2005-215051 20050222
CA 2556447 A1 20050901 CA 2005-2556447 20050222
EP 1723100 A2 20061122 EP 2005-713917 20050222
R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE,
IS, IT, LI, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, AL, BA,
HR, LV, MK, YU
US 2007123592 A1 20070531 US 2006-588478 20061002
PRAI US 2004-546111P P 20040219
WO 2005-US5557 W 20050222
OS CASREACT 143:248664; MARPAT 143:248664
GI



AB The invention relates to compds. and pharmaceutical compns. based on terephthalamide which are proteomimetic and methods for inhibiting the interaction of an α -helical protein with another protein or binding site and for treating diseases or conditions which are modulated through these interactions. Compds. I [X is H, halo, R, OR, SR or an amino group, where R is H, (un)substituted alkyl, acyl, aryl, heteroaryl, alkylenearyl or alkyleneheteroaryl; X₂, X₃, X₄ are independently H, halo, OH, Rc, ORc, where Rc is (un)substituted alkyl, acyl, aryl or alkylenearyl; R₄ is H, (un)substituted alkyl, alkenyl or alkylene amine; R', R_{1a}, R_{1b} are any group given for R₄ or (CH₂)₀₋₂CHR₂CO₂H or an alkyl ester, where R₂ is independently H or (un)substituted hydrocarbon, alkoxy, ester, alkanol, alkanoic acid, amine, etc.; or N-R' or NR_{1a}R_{1b} form an amino acid residue] are claimed. Thus, peptidomimetic compound II, prepared via coupling reaction of L-leucine Me ester hydrochloride, showed inhibitory constant K_i = 0.781 \pm 0.070 μ M in a fluorescence polarization assay (binding affinity for Bcl-XL).

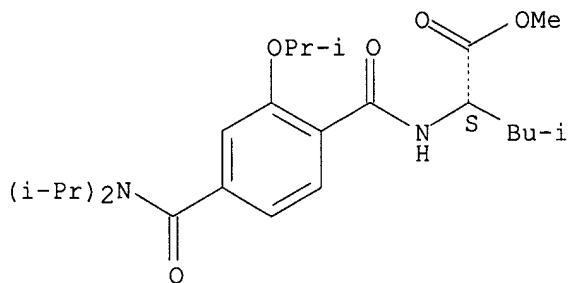
IT 681466-00-2P
RL: PAC (Pharmacological activity); PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(conformation; preparation of terephthalamide peptidomimetic compds. for therapeutic use)
RN 681466-00-2 CAPLUS
CN L-Leucine, N-[4-[[[(1S)-2-methoxy-1-methyl-2-oxoethyl](2-methylpropyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



IT 681465-54-3P
RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(preparation of terephthalamide peptidomimetic compds. for therapeutic use)
RN 681465-54-3 CAPLUS
CN L-Leucine, N-[4-[[bis(1-methylethyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



IT 681465-56-5P 681465-60-1P 681465-62-3P
 681465-64-5P 681465-66-7P 681465-68-9P
 681465-70-3P 681465-74-7P 852065-21-5P
 852065-22-6P 852065-23-7P 852065-25-9P
 852065-26-0P 852065-27-1P 852065-28-2P
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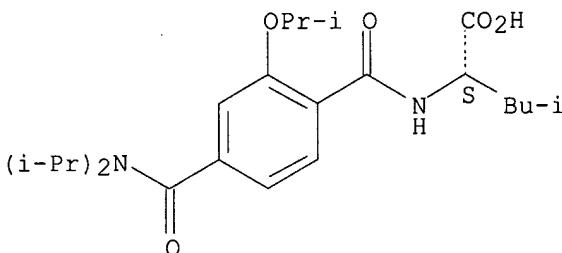
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of terephthalamide peptidomimetic compds. for therapeutic use)

RN 681465-56-5 CAPPLUS

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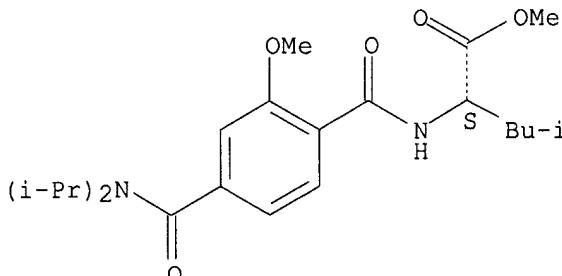
Absolute stereochemistry.



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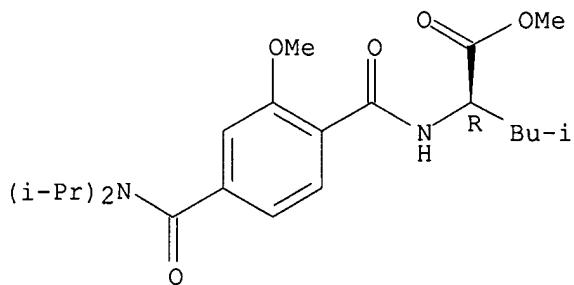
Absolute stereochemistry. Rotation (-).



RN 681465-62-3 CAPPLUS

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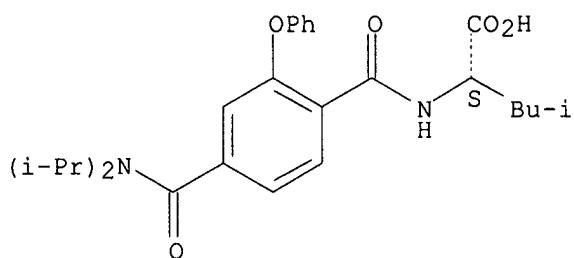
Absolute stereochemistry.



RN 681465-64-5 CAPLUS

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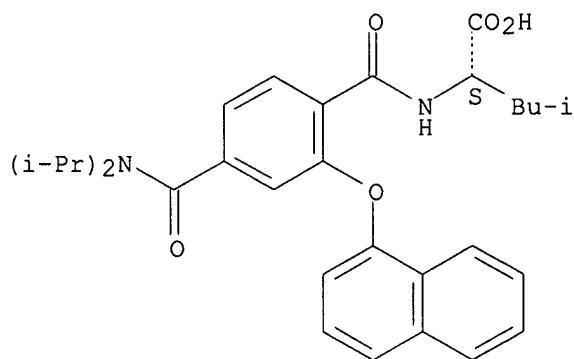
Absolute stereochemistry.



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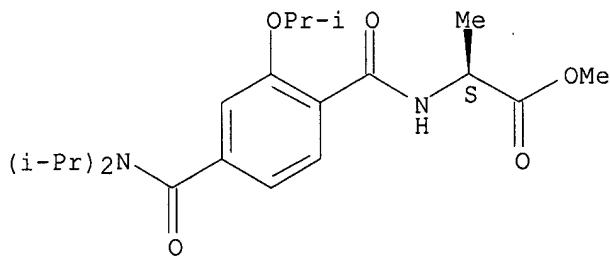
Absolute stereochemistry.



RN 681465-68-9 CAPLUS

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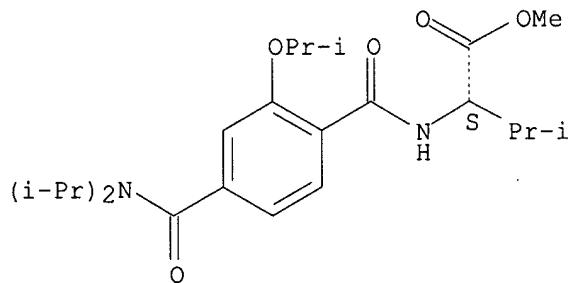
Absolute stereochemistry.



RN 681465-70-3 CAPLUS

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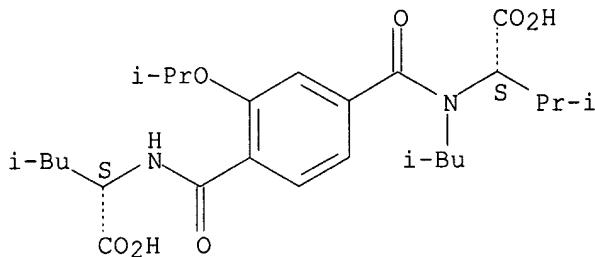
Absolute stereochemistry.



RN 681465-74-7 CAPLUS

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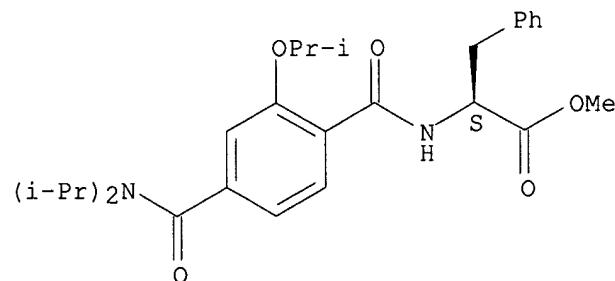
Absolute stereochemistry.



RN 852065-21-5 CAPLUS

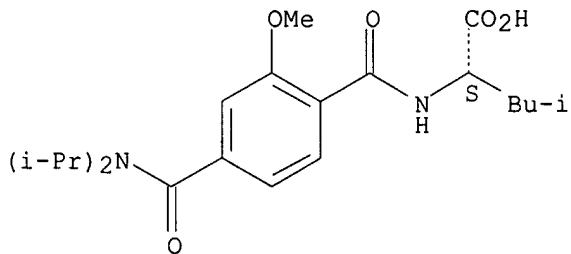
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Absolute stereochemistry.



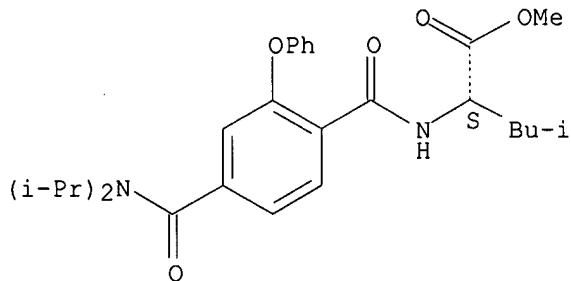
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(CA INDEX NAME)

Absolute stereochemistry.



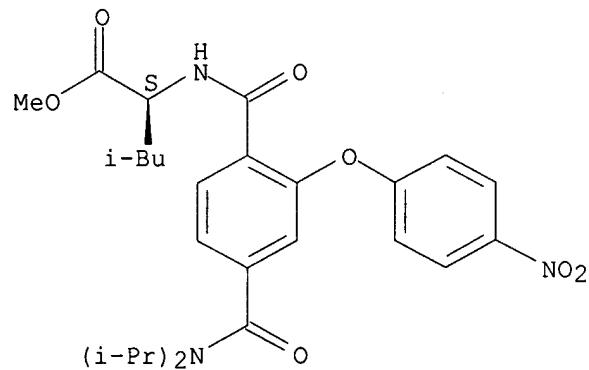
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Absolute stereochemistry.



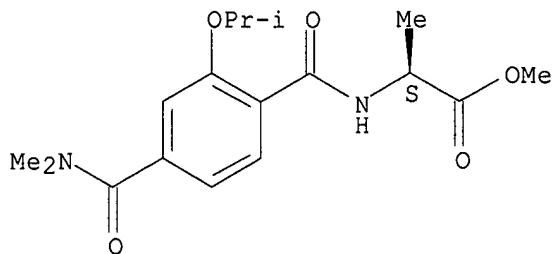
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Absolute stereochemistry.



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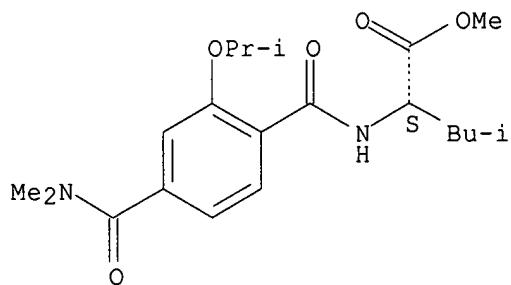
Absolute stereochemistry. Rotation (+).



RN 852065-27-1 CAPLUS

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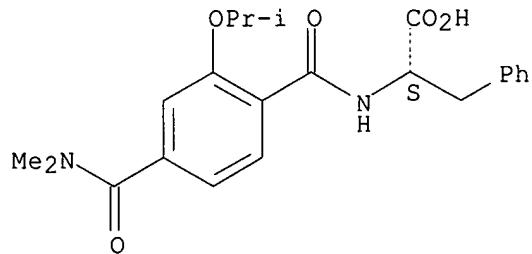
Absolute stereochemistry.



RN 852065-28-2 CAPLUS

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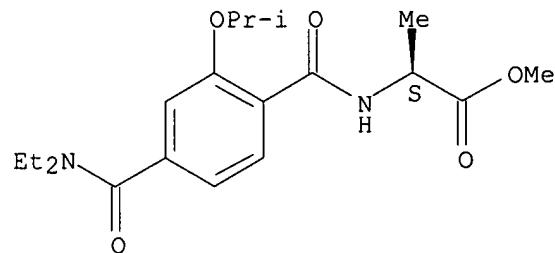
Absolute stereochemistry. Rotation (+).



RN 852065-29-3 CAPLUS

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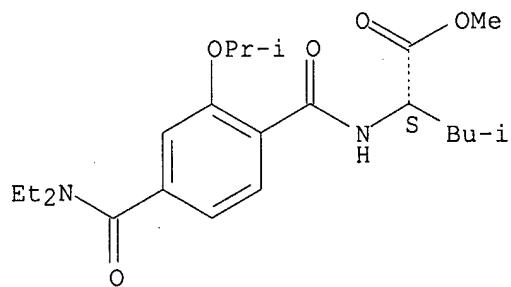
Absolute stereochemistry. Rotation (+).



RN 852065-30-6 CAPLUS

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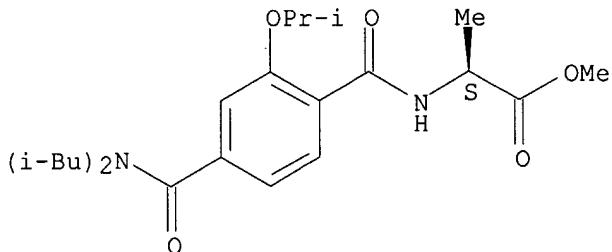
Absolute stereochemistry. Rotation (+).



RN 852065-31-7 CAPLUS

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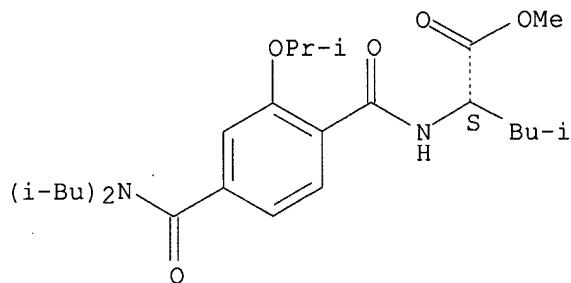
Absolute stereochemistry. Rotation (+).



RN 852065-32-8 CAPLUS

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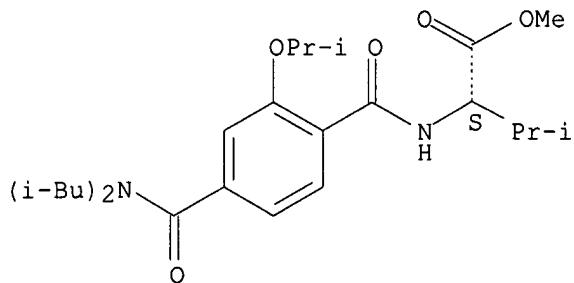
Absolute stereochemistry. Rotation (+).



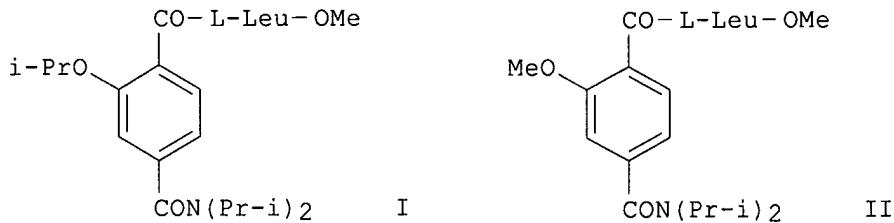
RN 852065-33-9 CAPLUS

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Absolute stereochemistry. Rotation (+).

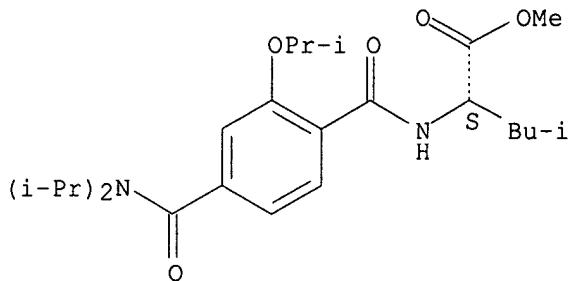


L7 ANSWER 2 OF 4 CAPLUS COPYRIGHT 2008 ACS on STN
 AN 2005:267006 CAPLUS
 DN 142:482288
 TI Terephthalamide Derivatives as Mimetics of Helical Peptides: Disruption of the Bcl-xL/Bak Interaction
 AU Yin, Hang; Lee, Gui-in; Sedey, Kristine A.; Rodriguez, Johanna M.; Wang, Hong-Gang; Sebti, Said M.; Hamilton, Andrew D.
 CS Department of Chemistry, Yale University, New Haven, CT, 06520-8107, USA
 SO Journal of the American Chemical Society (2005), 127(15), 5463-5468
 CODEN: JACSAT; ISSN: 0002-7863
 PB American Chemical Society
 DT Journal
 LA English
 OS CASREACT 142:482288
 GI



AB A series of Bcl-xL/Bak antagonists, based on a terephthalamide scaffold, was designed to mimic the α -helical region of the Bak peptide. These mols. showed favorable in-vitro activities in disrupting the Bcl-xL/Bak BH3 domain complex (terephthalamides I and II, $K_i = 0.78 \pm 0.07$ and $1.85 \pm 0.32 \mu\text{M}$, resp.). Extensive structure-affinity studies demonstrated a correlation between the ability of terephthalamide derivs. to disrupt Bcl-xL/Bak complex formation and the size of variable side chains on these mols. Treatment of human HEK293 cells with the terephthalamide derivative 26 resulted in disruption of the Bcl-xL/Bax interaction in whole cells with an IC_{50} of $35.0 \mu\text{M}$. Computational docking simulations and NMR expts. suggested that the binding cleft for the BH3 domain of the Bak peptide on the surface of Bcl-xL is the target area for these synthetic inhibitors.
 IT 681465-54-3P
 RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent)
 (preparation of terephthalamide derivs. as helical peptidomimetics and their evaluation as antagonists of Bcl-xL/Bak protein interaction)
 RN 681465-54-3 CAPLUS
 CN L-Leucine, N-[4-[[bis(1-methylethyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



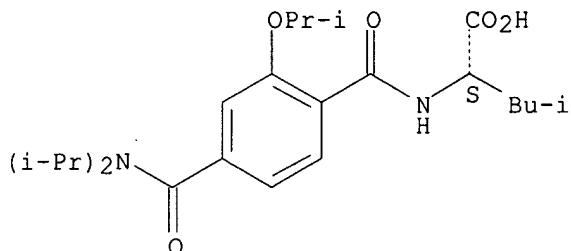
IT 681465-56-5P 681465-58-7P 681465-60-1P
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852065-28-2P 852065-29-3P 852065-30-6P
852065-31-7P 852065-32-8P 852065-33-9P

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); BIOL
(Biological study); PREP (Preparation)
(preparation of terephthalamide derivs. as helical peptidomimetics and their
evaluation as antagonists of Bcl-x/Bak protein interaction)

RN 681465-56-5 CAPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]- (CA INDEX NAME)

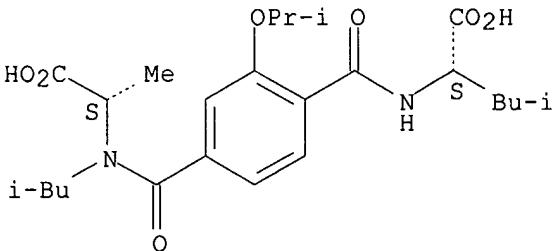
Absolute stereochemistry.



RN 681465-58-7 CAPLUS

CN L-Leucine, N-[4-[[[(1S)-1-carboxyethyl](2-methylpropyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]- (9CI) (CA INDEX NAME)

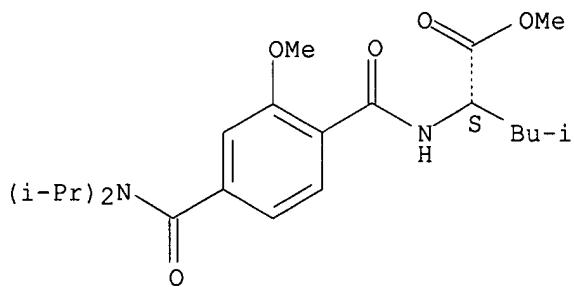
Absolute stereochemistry. Rotation (-).



RN 681465-60-1 CAPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-methoxybenzoyl]-, methyl ester (CA INDEX NAME)

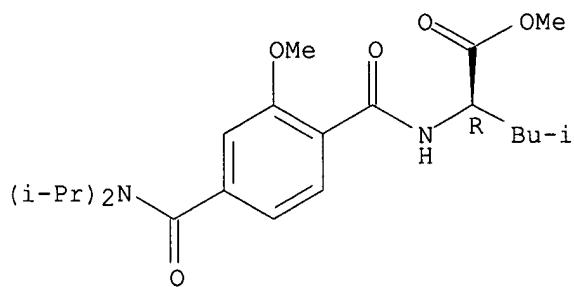
Absolute stereochemistry. Rotation (-).



RN 681465-62-3 CAPLUS

CN D-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-methoxybenzoyl]-, methyl ester (CA INDEX NAME)

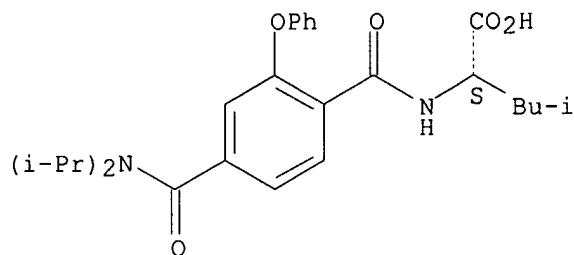
Absolute stereochemistry.



RN 681465-64-5 CAPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-phenoxybenzoyl]- (CA INDEX NAME)

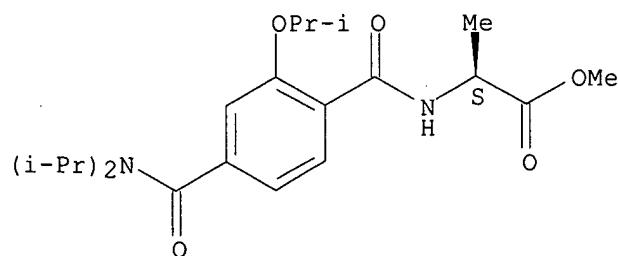
Absolute stereochemistry.



RN 681465-68-9 CAPLUS

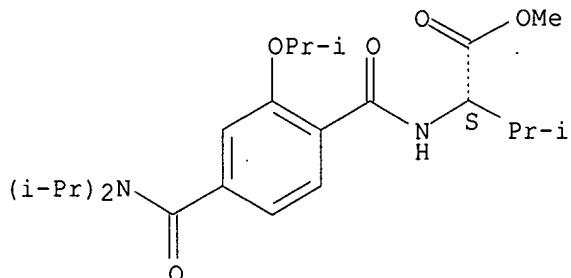
CN L-Alanine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

Absolute stereochemistry.



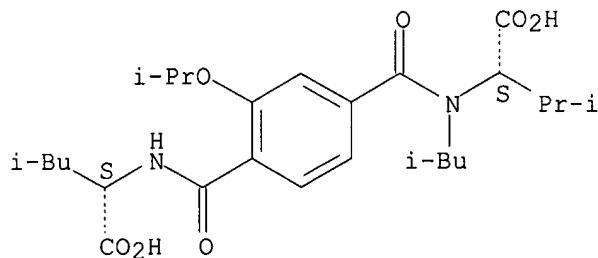
RN 681465-70-3 CAPLUS
CN L-Valine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

Absolute stereochemistry.



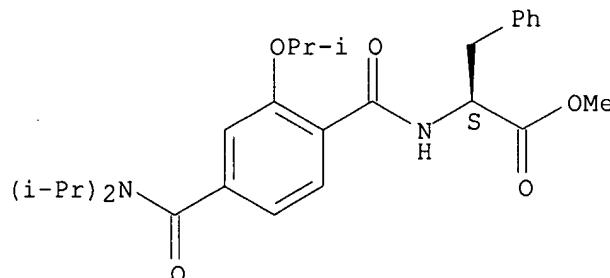
RN 681465-74-7 CAPLUS
CN L-Leucine, N-[4-[[[(1S)-1-carboxy-2-methylpropyl](2-methylpropyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



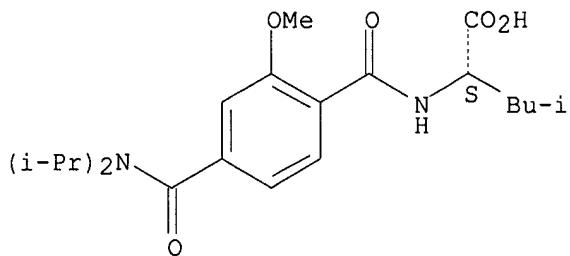
RN 852065-21-5 CAPLUS
CN L-Phenylalanine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

Absolute stereochemistry.



RN 852065-22-6 CAPLUS
CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-methoxybenzoyl]- (CA INDEX NAME)

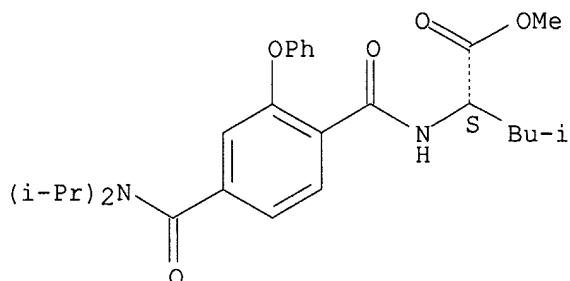
Absolute stereochemistry.



RN 852065-23-7 CAPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-phenoxybenzoyl]-, methyl ester (CA INDEX NAME)

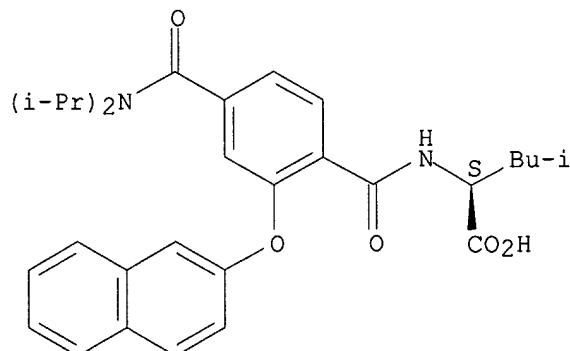
Absolute stereochemistry.



RN 852065-24-8 CAPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(2-naphthoxy)benzoyl]- (CA INDEX NAME)

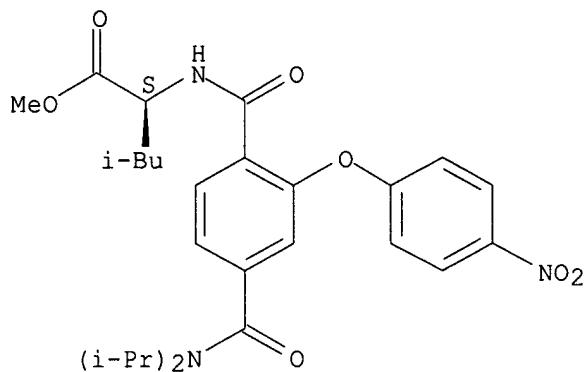
Absolute stereochemistry.



RN 852065-25-9 CAPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(4-nitrophenoxy)benzoyl]-, methyl ester (CA INDEX NAME)

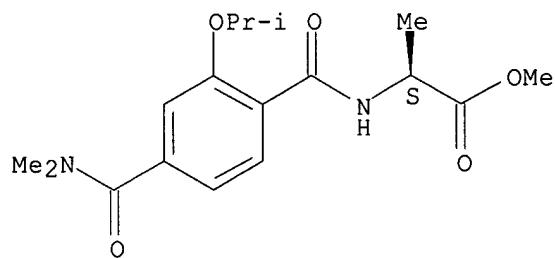
Absolute stereochemistry.



RN 852065-26-0 CAPLUS

CN L-Alanine, N-[4-[(dimethylamino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

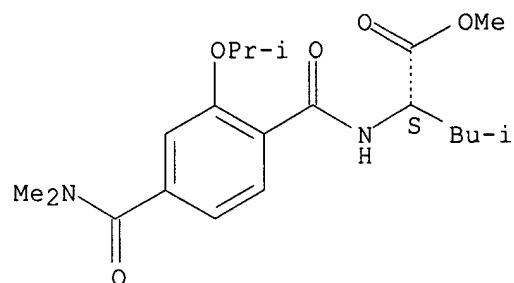
Absolute stereochemistry. Rotation (+).



RN 852065-27-1 CAPLUS

CN L-Leucine, N-[4-[(dimethylamino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

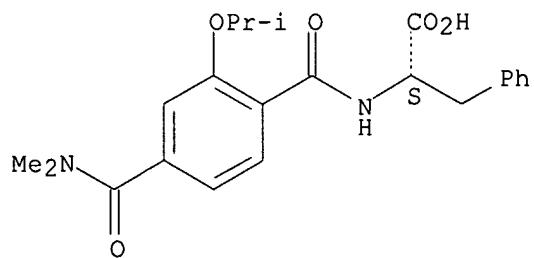
Absolute stereochemistry.



RN 852065-28-2 CAPLUS

CN L-Phenylalanine, N-[4-[(dimethylamino)carbonyl]-2-(1-methylethoxy)benzoyl]- (CA INDEX NAME)

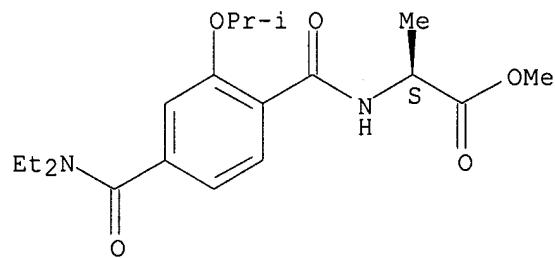
Absolute stereochemistry. Rotation (+).



RN 852065-29-3 CAPLUS

CN L-Alanine, N-[4-[(diethylamino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

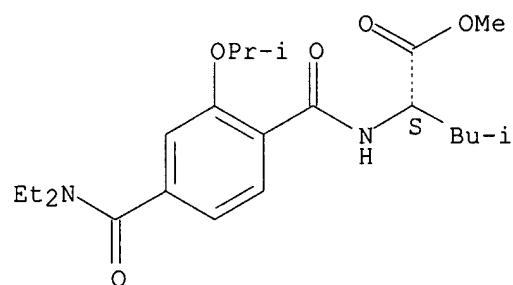
Absolute stereochemistry. Rotation (+).



RN 852065-30-6 CAPLUS

CN L-Leucine, N-[4-[(diethylamino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

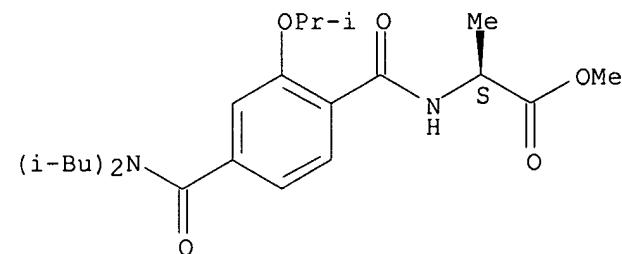
Absolute stereochemistry. Rotation (+).



RN 852065-31-7 CAPLUS

CN L-Alanine, N-[4-[(bis(2-methylpropyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

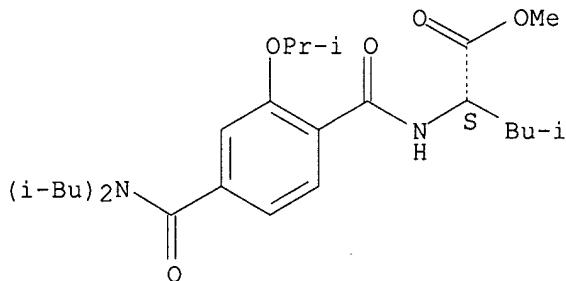
Absolute stereochemistry. Rotation (+).



RN 852065-32-8 CAPLUS

CN L-Leucine, N-[4-[(bis(2-methylpropyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

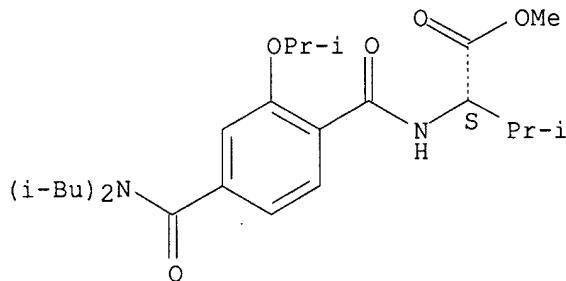
Absolute stereochemistry. Rotation (+).



RN 852065-33-9 CAPLUS

CN L-Valine, N-[4-[(bis(2-methylpropyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

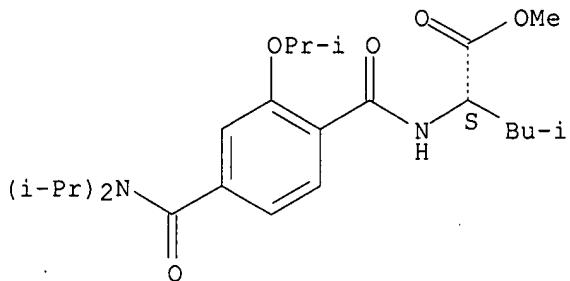
Absolute stereochemistry. Rotation (+).



RE.CNT 28 THERE ARE 28 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 3 OF 4 CAPLUS COPYRIGHT 2008 ACS on STN
AN 2004:189140 CAPLUS
DN 140:350050
TI Terephthalamide derivatives as mimetics of the helical region of Bak peptide target Bcl-xL protein
AU Yin, Hang; Hamilton, Andrew D.
CS Department of Chemistry, Yale University, New Haven, CT, 06511, USA
SO Bioorganic & Medicinal Chemistry Letters (2004), 14(6), 1375-1379
CODEN: BMCLE8; ISSN: 0960-894X
PB Elsevier Science B.V.
DT Journal
LA English
OS CASREACT 140:350050
AB A group of novel Bcl-xL/Bak antagonists, based on a terephthalamide scaffold, were designed to mimic the α -helical region of the Bak peptide. Good in vitro inhibition potencies in disrupting the Bak/Bcl-xL complex have been observed (terephthalamide 4, $K_i=0.78\pm0.07$ μ M).
IT 681465-54-3P
RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(terephthalamide derivs. as novel Bcl-xL/Bak antagonists)
RN 681465-54-3 CAPLUS
CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



IT 681465-56-5P 681465-58-7P

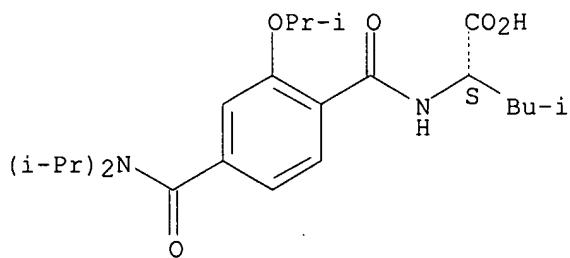
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(terephthalamide derivs. as novel Bcl-xL/Bak antagonists)

RN 681465-56-5 CAPPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(1-methylethoxy)benzoyl]- (CA INDEX NAME)

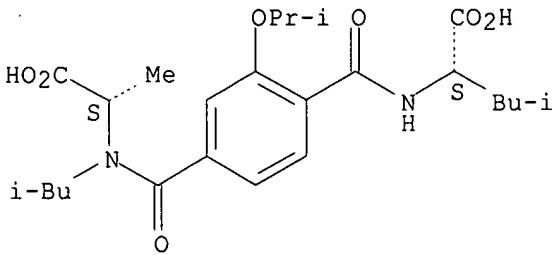
Absolute stereochemistry.



RN 681465-58-7 CAPPLUS

CN L-Leucine, N-[4-[[[(1S)-1-carboxyethyl](2-methylpropyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



IT 681465-60-1 681465-62-3 681465-64-5

681465-66-7 681465-68-9 681465-70-3

681465-72-5 681465-74-7

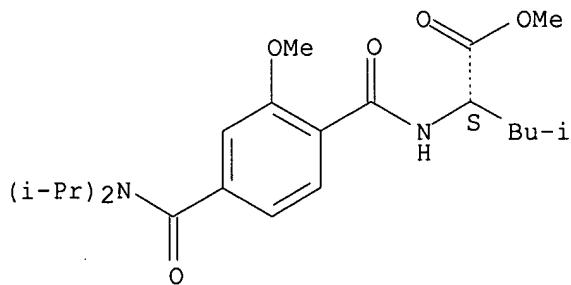
RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

(terephthalamide derivs. as novel Bcl-xL/Bak antagonists)

RN 681465-60-1 CAPPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-methoxybenzoyl]-, methyl ester (CA INDEX NAME)

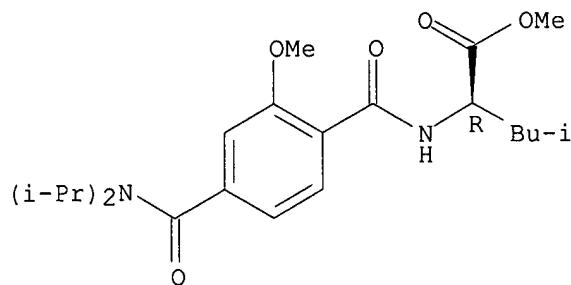
Absolute stereochemistry. Rotation (-).



RN 681465-62-3 CAPLUS

CN D-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-methoxybenzoyl]-, methyl ester (CA INDEX NAME)

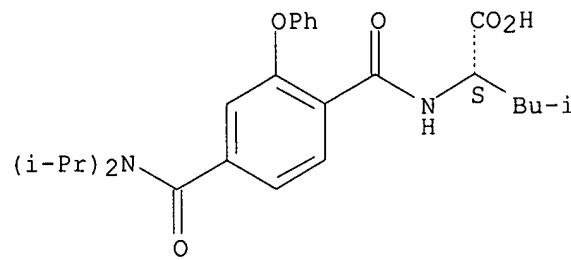
Absolute stereochemistry.



RN 681465-64-5 CAPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-phenoxybenzoyl]- (CA INDEX NAME)

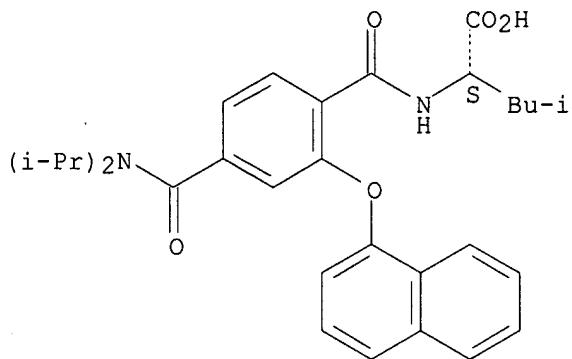
Absolute stereochemistry.



RN 681465-66-7 CAPLUS

CN L-Leucine, N-[4-[(bis(1-methylethyl)amino)carbonyl]-2-(1-naphthalenyl)benzoyl]- (CA INDEX NAME)

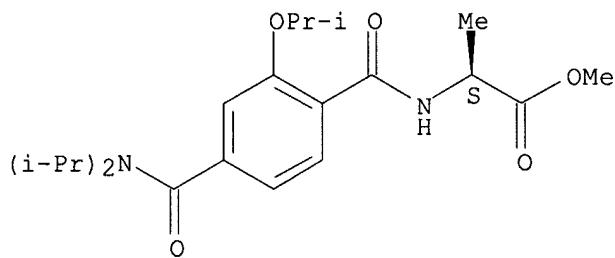
Absolute stereochemistry.



RN 681465-68-9 CAPLUS

CN L-Alanine, N-[4-[[bis(1-methylethyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

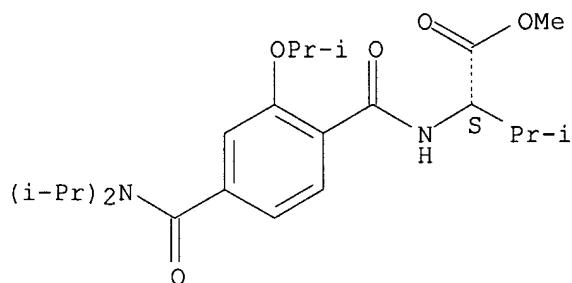
Absolute stereochemistry.



RN 681465-70-3 CAPLUS

CN L-Valine, N-[4-[[bis(1-methylethyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (CA INDEX NAME)

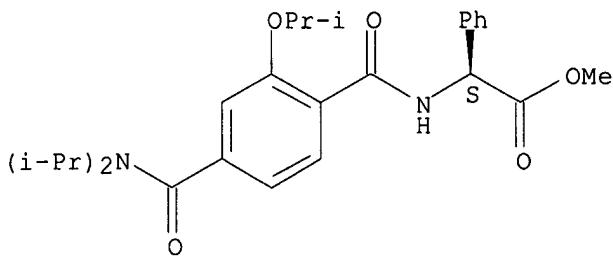
Absolute stereochemistry.



RN 681465-72-5 CAPLUS

CN Benzeneacetic acid, α -[[4-[[bis(1-methylethyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]amino]-, methyl ester, (α S)- (CA INDEX NAME)

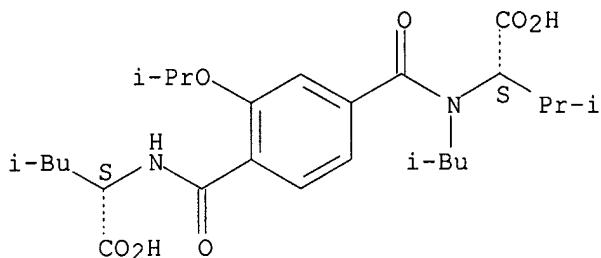
Absolute stereochemistry.



RN 681465-74-7 CAPLUS

CN L-Leucine, N-[4-[[[(1S)-1-carboxy-2-methylpropyl](2-methylpropyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



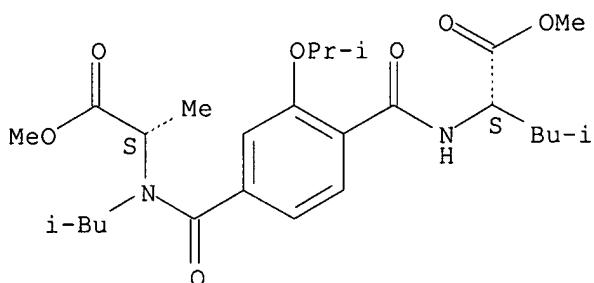
IT 681466-00-2P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(terephthalamide derivs. as novel Bcl-xL/Bak antagonists)

RN 681466-00-2 CAPLUS

CN L-Leucine, N-[4-[[[(1S)-2-methoxy-1-methyl-2-oxoethyl](2-methylpropyl)amino]carbonyl]-2-(1-methylethoxy)benzoyl]-, methyl ester (9CI) (CA INDEX NAME)

Absolute stereochemistry. Rotation (-).



RE.CNT 15 THERE ARE 15 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L7 ANSWER 4 OF 4 CAPLUS COPYRIGHT 2008 ACS on STN

AN 1998:394328 CAPLUS

DN 129:67773

TI Preparation of benzamide derivatives having a vasopressin antagonistic activity

IN Setoi, Hiroyuki; Ohkawa, Takehiko; Zenkoh, Tatsuya; Sawada, Hitoshi; Sawada, Yuki; Oku, Teruo

PA Fujisawa Pharmaceutical Co., Ltd., Japan

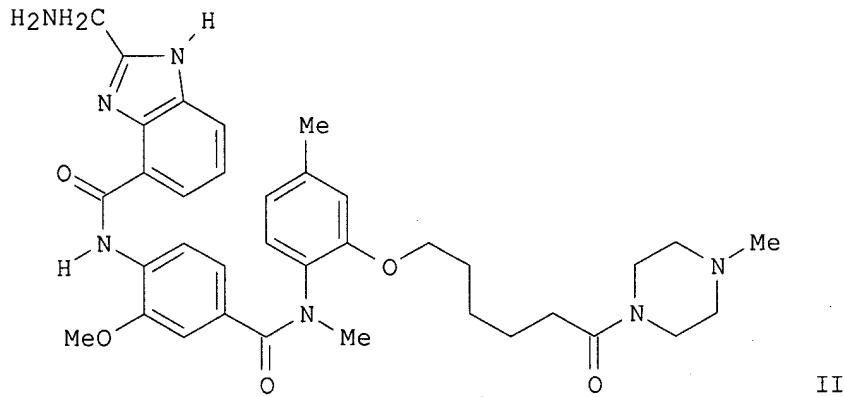
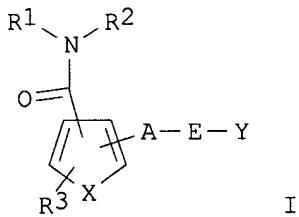
SO PCT Int. Appl., 332 pp.

CODEN: PIXXD2

DT Patent
LA English

FAN.CNT 1

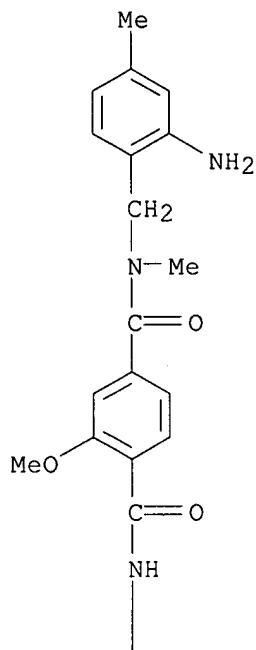
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PI	WO 9824771	A1	19980611	WO 1997-JP4192	19971118
	W: AU, CA, CN, HU, IL, JP, KR, MX, US, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE				
AU	9749672	A	19980629	AU 1997-49672	19971118
EP	946519	A1	19991006	EP 1997-912493	19971118
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, PT, IE, FI				
JP	2001505193	T	20010417	JP 1998-521225	19971118
US	6207693	B1	20010327	US 1999-308662	19990602
	US 6316482	B1	20011113	US 2000-614132	20000711
PRAI	AU 1996-3953	A	19961202		
	WO 1997-JP4192	W	19971118		
	US 1999-308662	A3	19990602		
OS	MARPAT 129:67773				
GI					



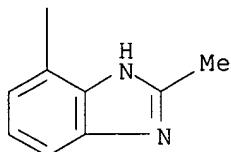
AB The title compds. [I; R1 = (un)substituted aryl, cyclo(lower)alkyl, heterocyclyl; R2 = H, lower alkyl, etc.; R3 = H, halo, OH, etc.; A = a single bond, O, NH; E = lower alkylene, lower alkenylene, etc.; X = CH:CH, CH:N, S; Y = (un)substituted aryl, condensed heterocyclyl, etc.] and their pharmaceutically acceptable salts, useful in treatment and/or prevention of hypertension, heart failure, renal insufficiency, edema, ascites, vasopressin parasecretion syndrome, hepatocirrhosis, hyponatremia, hypokalemia, diabetic, circulation disorder, cerebrovascular disease, Meniere's disease or motion sickness, were prepared. Thus, the title compound II showed IC50 of 1.5 nM against vasopressin 1 receptor binding.

IT 208770-26-7P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of benzamide derivs. having a vasopressin antagonistic activity)
RN 208770-26-7 CAPLUS
CN 1,4-Benzenedicarboxamide, N4-[(2-amino-4-methylphenyl)methyl]-2-methoxy-N4-methyl-N1-(2-methyl-1H-benzimidazol-4-yl)-, dihydrochloride (9CI) (CA INDEX NAME)

PAGE 1-A



PAGE 2-A



● 2 HCl

RE.CNT 17 THERE ARE 17 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT